

By: Senator(s) Tate

To: Drug Policy

SENATE BILL NO. 2110

1 AN ACT TO AMEND SECTION 41-29-113, MISSISSIPPI CODE OF 1972,
2 TO ADD KRATOM TO SCHEDULE I OF THE UNIFORM CONTROLLED SUBSTANCES
3 ACT; AND FOR RELATED PURPOSES.

4 BE IT ENACTED BY THE LEGISLATURE OF THE STATE OF MISSISSIPPI:

5 **SECTION 1.** Section 41-29-113, Mississippi Code of 1972, is
6 amended as follows:

7 41-29-113.

8 **SCHEDULE I**

9 (a) Schedule I consists of the drugs and other substances,
10 by whatever official name, common or usual name, chemical name, or
11 brand name designated, that is listed in this section.

12 (b) **Opiates.** Unless specifically excepted or unless listed
13 in another schedule, any of the following opiates, including their
14 isomers, esters, ethers, salts and salts of isomers, esters and
15 ethers, whenever the existence of these isomers, esters, ethers
16 and salts is possible within the specific chemical designation:

17 (1) Acetyl-alpha-methylfentanyl;



- 18 (2) Acetyl Fentanyl
19 N-(1-phenethylpiperidin-4-yl)-N-phenylacetamide;
20 (3) AH-7921 (3,4-dichloro-N-[(1-dimethylamino)
21 cyclohexylmethyl]benzamide);
22 (4) Acetylmethadol;
23 (5) Allylprodine;
24 (6) Alphacetylmethadol, except levo-alphacetylmethadol
25 (levo-alpha-acetylmethadol, levomethadyl acetate, or LAAM);
26 (7) Alphameprodine;
27 (8) Alphamethadol;
28 (9) Alpha-methylfentanyl;
29 (10) Alpha-methylthiofentanyl;
30 (11) Benzethidine;
31 (12) Betacetylmethadol;
32 (13) Beta-hydroxyfentanyl;
33 (14) Beta-hydroxy-3-methylfentanyl;
34 (15) Betameprodine;
35 (16) Betamethadol;
36 (17) Betaprodine;
37 (18) Butyrl fentanyl
38 (N-(1-phenethylpiperidin-4-yl)-N-phenylbutyramide);
39 (19) Clonitazene;
40 (20) Dextromoramide;
41 (21) Diampromide;
42 (22) Diethylthiambutene;



- 43 (23) DifenoXin;
- 44 (24) Dimenoxadol;
- 45 (25) Dimepheptanol;
- 46 (26) Dimethylthiambutene;
- 47 (27) Dioxaphetyl butyrate;
- 48 (28) Dipipanone;
- 49 (29) Ethylmethylthiambutene;
- 50 (30) Etonitazene;
- 51 (31) EtoXeridine;
- 52 (32) Fentanyl-related substances, meaning any substance
- 53 not otherwise listed under another schedule and for which no
- 54 exemption or approval is in effect under Section 505 of the
- 55 Federal Food, Drug, and Cosmetic Act [21 USC 355] that is
- 56 structurally related to fentanyl by one or more of the following
- 57 modifications:
- 58 (A) Replacement of the phenyl portion of the
- 59 phenethyl group by any monocycle, whether or not further
- 60 substituted in or on the monocycle;
- 61 (B) Substitution in or on the phenethyl group with
- 62 alkyl, alkenyl, alkoXyl, hydroXyl, halo, haloalkyl, amino or nitro
- 63 groups;
- 64 (C) Substitution in or on the piperidine ring with
- 65 alkyl, alkenyl, alkoXyl, ester, ether, hydroXyl, halo, haloalkyl,
- 66 amino or nitro groups;



67 (D) Replacement of the aniline ring with any
68 aromatic monocycle whether or not further substituted in or on the
69 aromatic monocycle; and/or

70 (E) Replacement of the N-propionyl group by
71 another acyl group.

72 Fentanyl-related substances include, but are not limited to,
73 cyclopropyl fentanyl,

74 (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopropanecarboxamide);

75 Furanyl-Fentanyl,

76 (N-(1-phenethylpiperidin-4-yl)-N-phenylfuran-2-carboxamide);

77 valeryl fentanyl,

78 (N-(1-phenethylpiperidin-4-yl)-N-phenylpentanamide);

79 para-fluorobutyryl fentanyl,

80 (N-(4-fluorophenyl)-N-(1-phenethylpiperidin-4-yl)butyramide);

81 para-methoxybutyryl fentanyl,

82 (N-(4-methoxyphenyl)-N-(1-phenethylpiperidin-4-yl)butyramide);

83 para-chloroisobutyryl fentanyl,

84 (N-(4-chlorophenyl)-N-(1-phenethylpiperidin-4-yl)isobutyramide);

85 isobutyryl fentanyl,

86 (N-(1-phenethylpiperidin-4-yl)-N-phenylisobutyramide);

87 cyclopentyl fentanyl,

88 (N-(1-phenethylpiperidin-4-yl)-N-phenylcyclopentanecarboxamide);

89 and



90 ocfentanil,
91 (N-(2-fluorophenyl)-2-methoxy-N-(1-phenethylpiperidin-4-yl)acetami
92 de);
93 (33) Furethidine;
94 (34) Hydroxypethidine;
95 (35) Ketobemidone (including the optical and geometric
96 isomers);
97 (36) Levomoramide;
98 (37) Levophenacilmorphan;
99 (38) 3-methylfentanyl;
100 (39) 3-methylthiofentanyl;
101 (40) Morpheridine;
102 (41) MPPP (1-methyl-4-phenyl-4-propionoxypiperidine);
103 (42)
104 *N*-[1-[2-hydroxy-2-(thiophen-2-yl)ethyl]piperidin-4-yl]-*N*-phenylprop
105 ionamide, its isomers, esters, ethers, salts and salts of isomers,
106 esters and ethers (other names: beta-hydroxythiofentanyl);
107 (43) Noracymethadol;
108 (44) Norlevorphanol;
109 (45) Normethadone;
110 (46) Norpipanone;
111 (47) Para-fluorofentanyl;
112 (48) PEPAP
113 (1-(-2-phenethyl)-4-phenyl-4-acetoxypiperidine);
114 (49) Phenadoxone;



- 115 (50) Phenampromide;
116 (51) Phenomorphan;
117 (52) Phenoperidine;
118 (53) Piritramide;
119 (54) Proheptazine;
120 (55) Properidine;
121 (56) Propiram;
122 (57) Racemoramide;
123 (58) Thiofentanyl;
124 (59) Tilidine;
125 (60) Trimeperidine;
126 (61) U-47700,
127 3,4-dichloro-N-[2-(dimethylamino)cyclohexyl]-N-methylbenzamide.

128 (c) **Opium derivatives.** Unless specifically excepted or
129 unless listed in another schedule, any of the following opium
130 derivatives, their salts, isomers and salts of isomers, whenever
131 the existence of these salts, isomers and salts of isomers is
132 possible within the specific chemical designation:

- 133 (1) Acetorphine;
134 (2) Acetyldihydrocodeine;
135 (3) Benzylmorphine;
136 (4) Codeine methylbromide;
137 (5) Codeine-N-Oxide;
138 (6) Cyrenorphine;
139 (7) Desomorphine;



- 140 (8) Dihydromorphine;
- 141 (9) Drotebanol;
- 142 (10) Etorphine (except hydrochloride salt);
- 143 (11) Heroin;
- 144 (12) Hydromorphinol;
- 145 (13) Methyldesorphine;
- 146 (14) Methyldihydromorphine;
- 147 (15) Monoacetylmorphine;
- 148 (16) Morphine methylbromide;
- 149 (17) Morphine methylsulfonate;
- 150 (18) Morphine-N-Oxide;
- 151 (19) Myrophine;
- 152 (20) Nicocodeine;
- 153 (21) Nicomorphine;
- 154 (22) Normorphine;
- 155 (23) Pholcodine;
- 156 (24) Thebacon.

157 (d) **Hallucinogenic substances.** Unless specifically excepted
158 or unless listed in another schedule, any material, compound,
159 mixture or preparation which contains any quantity of the
160 following substances, their salts, isomers (whether optical,
161 positional, or geometric) and salts of isomers, whenever the
162 existence of these salts, isomers and salts of isomers is possible
163 within the specific chemical designation:

- 164 (1) Alpha-ethyltryptamine;



- 165 (2) 4-bromo-2,5-dimethoxy-amphetamine;
- 166 (3) 4-bromo-2,5-dimethoxyphenethylamine;
- 167 (4) 2,5-dimethoxyamphetamine;
- 168 (5) 2,5-dimethoxy-4-ethylamphetamine (DOET);
- 169 (6) 2,5-dimethoxy-4-(n)-propylthiophenethylamine
170 (2C-T-7);
- 171 (7) 4-methoxyamphetamine;
- 172 (8) 5-methoxy-3,4-methylenedioxy-amphetamine;
- 173 (9) 4-methyl-2,5-dimethoxy-amphetamine;
- 174 (10) 3,4-methylenedioxy amphetamine;
- 175 (11) 3,4-methylenedioxymethamphetamine (MDMA);
- 176 (12) 3,4-methylenedioxy-N-ethylamphetamine (also known
177 as N-ethyl-alpha-methyl-3,4(methylenedioxy)phenethylamine, N-ethyl
178 MDA, MDE, MDEA);
- 179 (13) N-hydroxy-3,4-methylenedioxyamphetamine (also
180 known as N-hydroxy MDA, N-OHMDA, and
181 N-hydroxy-alpha-methyl-3,4(methylenedioxy)phenethylamine);
- 182 (14) 3,4,5-trimethoxy amphetamine;
- 183 (15) 5-methoxy-N,N-dimethyltryptamine (5-MeO-DMT);
- 184 (16) Alpha-methyltryptamine (also known as AMT);
- 185 (17) Bufotenine;
- 186 (18) Diethyltryptamine;
- 187 (19) Dimethyltryptamine;
- 188 (20) 5-methoxy-N,N-diisopropyltryptamine (5-MeO-DIPT);
- 189 (21) Ibogaine;



- 190 (22) Lysergic acid diethylamide (LSD);
- 191 (23) (A) Marijuana (Hemp as defined and regulated
192 under Sections 69-25-201 through 69-25-221 and Cannabidiol
193 contained in a legend drug product approved by the Federal Food
194 and Drug Administration or obtained under Section 41-29-136 are
195 exempt under Schedule I);
- 196 (B) Hashish;
- 197 (24) Mescaline;
- 198 (25) Parahexyl;
- 199 (26) Peyote;
- 200 (27) N-ethyl-3-piperidyl benzilate;
- 201 (28) N-methyl-3-piperidyl benzilate;
- 202 (29) Psilocybin;
- 203 (30) Psilocyn;
- 204 (31) Tetrahydrocannabinols, meaning
205 tetrahydrocannabinols contained in a plant of the genus Cannabis
206 (cannabis plant), as well as the synthetic equivalents of the
207 substances contained in the cannabis plant, or in the resinous
208 extractives of such plant, and/or synthetic substances,
209 derivatives, and their isomers with similar chemical structure and
210 pharmacological activity to those substances contained in the
211 plant such as the following:
- 212 (A) 1 cis or trans tetrahydrocannabinol;
- 213 (B) 6 cis or trans tetrahydrocannabinol;
- 214 (C) 3,4 cis or trans tetrahydrocannabinol.



215 (Since nomenclature of these substances is not
216 internationally standardized, compounds of these structures,
217 regardless of atomic positions, are covered.)

218 ("Tetrahydrocannabinols" excludes dronabinol and nabilone.)
219 For purposes of this paragraph, tetrahydrocannabinols do not
220 include hemp or hemp products regulated under Sections 69-25-201
221 through 69-25-221.

222 However, the following products are exempted from control:

223 (i) THC-containing industrial products made
224 from cannabis stalks (e.g., paper, rope and clothing);

225 (ii) Processed cannabis plant materials used
226 for industrial purposes, such as fiber retted from cannabis stalks
227 for use in manufacturing textiles or rope;

228 (iii) Animal feed mixtures that contain
229 sterilized cannabis seeds and other ingredients (not derived from
230 the cannabis plant) in a formula designed, marketed and
231 distributed for nonhuman consumption;

232 (iv) Personal care products that contain oil
233 from sterilized cannabis seeds, such as shampoos, soaps, and body
234 lotions (if the products do not cause THC to enter the human
235 body);

236 (v) Hemp as regulated under Sections
237 69-25-201 through 69-25-221; and



238 (vi) Any product derived from the hemp plant
239 designed for human ingestion and/or consumption that is approved
240 by the United States Food and Drug Administration;

241 (32) Phencyclidine;

242 (33) Ethylamine analog of phencyclidine (PCE);

243 (34) Pyrrolidine analog of phencyclidine (PHP, PCPy);

244 (35) Thiophene analog of phencyclidine;

245 (36) 1-[1-(2-thienyl)cyclohexyl] pyrrolidine (TCPy);

246 (37) 4-methylmethcathinone (mephedrone);

247 (38) 3,4-methylenedioxypropylvalerone (MDPV);

248 (39) 2-(2,5-dimethoxy-4-ethylphenyl)ethanamine (2C-E);

249 (40) 2-(2,5-dimethoxy-4-methylphenyl)ethanamine (2C-D);

250 (41) 2-(4-chloro-2,5-dimethoxyphenyl)ethanamine (2C-C);

251 (42) 2-(4-iodo-2,5-dimethoxyphenyl)ethanamine (2C-I);

252 or 2,5-dimethoxy-4-iodophenethylamine;

253 (43) 2-[4-(ethylthio)-2,5-dimethoxyphenyl]ethanamine

254 (2C-T-2);

255 (44)

256 2-[4-(isopropylthio)-2,5-dimethoxyphenyl]ethanamine (2C-T-4);

257 (45) 2-(2,5-dimethoxyphenyl)ethanamine (2C-H);

258 (46) 2-(2,5-dimethoxy-4-nitro-phenyl)ethanamine (2C-N);

259 (47) 2-(2,5-dimethoxy-4-(n)-propylphenyl)ethanamine

260 (2C-P);

261 (48) 3,4-methylenedioxy-N-methylcathinone (methylone);



262 (49)
263 2-(4-bromo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine
264 (25B-NBOMe; 2C-B-NBOMe; 25B; Cimbi-36);
265 (50)
266 2-(4-chloro-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine
267 (25C-NBOMe; 2C-C-NBOMe; 25C; Cimbi-82);
268 (51)
269 2-(4-iodo-2,5-dimethoxyphenyl)-N-(2-methoxybenzyl)ethanamine or
270 N-[(2-methoxyphenyl)methyl]ethanamine (25I-NBOMe; 2C-I-NBOMe; 25I;
271 Cimbi-5);
272 (52) 7-bromo-5-(2-chlorophenyl)-1,3-dihydro-2H-1,
273 4-benzodiazepin-2-one (also known as Phenazepam);
274 (53) 7-(2-chlorophenyl)-4-ethyl-13-methyl-3-thia-1,8,
275 11,12-tetraazatricyclo[8.3.0.0]trideca-2(6),4,7,10,12-pentaene
276 (also known as Etizolam);
277 (54) Salvia divinorum;
278 (55) Synthetic cannabinoids. Unless specifically
279 excepted or unless listed in another schedule, any material,
280 compound, mixture, or preparation which contains any quantity of a
281 synthetic cannabinoid found in any of the following chemical
282 groups, whether or not substituted to any extent, or any of those
283 groups which contain any synthetic cannabinoid salts, isomers, or
284 salts of isomers, whenever the existence of such salts, isomers,
285 or salts of isomers is possible within the specific chemical



286 designation, including all synthetic cannabinoid chemical
287 analogues in such groups:

288 (A) (6aR,10aR)-9-(hydroxymethyl)-6,
289 6-dimethyl-3-(2-methyloctan-2-yl)-6a,7,10,10a-tetrahydrobenzo[c]
290 chromen-1-ol (also known as HU-210 or
291 1,1-dimethylheptyl-11-hydroxy-delta8-tetrahydrocannabinol);

292 (B) Naphthoylindoles and naphthylmethylinindoles,
293 being any compound structurally derived from 3-(1-naphthoyl)indole
294 or 1H-indol-3-yl-(1-naphthyl)methane, whether or not substituted
295 in the indole ring to any extent, or in the naphthyl ring to any
296 extent;

297 (C) Naphthoylpyrroles, being any compound
298 structurally derived from 3-(1-naphthoyl)pyrrole, whether or not
299 substituted in the pyrrole ring to any extent, or in the naphthyl
300 ring to any extent;

301 (D) Naphthylmethylindenes, being any compound
302 structurally derived from 1-(1-naphthylmethyl)indene, whether or
303 not substituted in the indene ring to any extent or in the
304 naphthyl ring to any extent;

305 (E) Phenylacetylindoles, being any compound
306 structurally derived from 3-phenylacetylindole, whether or not
307 substituted in the indole ring to any extent or in the phenyl ring
308 to any extent;

309 (F) Cyclohexylphenols, being any compound
310 structurally derived from 2-(3-hydroxycyclohexyl)phenol, whether



311 or not substituted in the cyclohexyl ring to any extent or in the
312 phenolic ring to any extent;

313 (G) Benzoylindoles, whether or not substituted in
314 the indole ring to any extent or in the phenyl ring to any extent;

315 (H) Adamantoylindoles, whether or not substituted
316 in the indole ring to any extent or in the adamantoyl ring system
317 to any extent;

318 (I) Tetrahydro derivatives of cannabinal and
319 3-alkyl homologues of cannabinal or of its tetrahydro derivatives,
320 except where contained in cannabis or cannabis resin;

321 (J) 3-Cyclopropylmethanone indole or
322 3-Cyclobutylmethanone indole or 3-Cyclopentylmethanone indole by
323 substitution at the nitrogen atom of the indole ring, whether or
324 not further substituted in the indole ring to any extent, whether
325 or not substituted on the cyclopropyl, cyclobutyl or cyclopentyl
326 rings to any extent;

327 (K) Quinoliny ester indoles, being any compound
328 structurally derived from 1H-indole-3carboxylic acid-8-quinoliny
329 ester, whether or not substituted in the indole ring to any extent
330 or the quinolone ring to any extent;

331 (L) 3-carboxamide-1H-indazoles, whether or not
332 substituted in the indazole ring to any extent and substituted to
333 any degree on the carboxamide nitrogen and
334 3-carboxamide-1H-indoles, whether or not substituted in the indole



335 ring to any extent and substituted to any degree on the
336 carboxamide nitrogen;

337 (M) Cycloalkanemethanone Indoles, whether or not
338 substituted at the nitrogen atom on the indole ring, whether or
339 not further substituted in the indole ring to any extent, whether
340 or not substituted on the cycloalkane ring to any extent.

341 (e) **Depressants.** Unless specifically excepted or unless
342 listed in another schedule, any material, compound, mixture, or
343 preparation which contains any quantity of the following
344 substances having a depressant effect on the central nervous
345 system, including their salts, isomers, and salts of isomers,
346 whenever the existence of such salts, isomers, and salts of
347 isomers is possible within the specific chemical designation:

348 (1) Gamma-hydroxybutyric acid (other names include:
349 GHB, gamma-hydroxybutyrate; 4-hydroxybutyrate; 4-hydroxybutanoic
350 acid; sodium oxybate; sodium oxybutyrate);

351 (2) Mecloqualone;

352 (3) Methaqualone.

353 (f) **Stimulants.** Any material, compound, mixture or
354 preparation which contains any quantity of the following central
355 nervous system stimulants including optical salts, isomers and
356 salts of isomers unless specifically excepted or unless listed in
357 another schedule:

358 (1) Aminorex;



359 (2) N-benzylpiperazine (also known as BZP and
360 1-benzylpiperazine);

361 (3) Cathinone;

362 (4) Fenethylamine;

363 (5) Methcathinone;

364 (6) 4-methylaminorex (also known as
365 2-amino-4-methyl-5-phenyl-2-oxazoline);

366 (7) N-ethylamphetamine;

367 (8) Any material, compound, mixture or preparation
368 which contains any quantity of N,N-dimethylamphetamine. (Other
369 names include: N,N,-alpha-trimethyl-benzeneethanamine and
370 N,N-alpha-trimethylphenethylamine);

371 (9) **Synthetic cathinones.** (A) Unless listed in
372 another schedule, any compound other than bupropion that is
373 structurally derived from 2-Amino-1-phenyl-1-propanone by
374 modification in any of the following ways:

375 (i) By substitution in the phenyl ring to any
376 extent with alkyl, alkoxy, alkylendioxy, haloalkyl or halide
377 substituents, whether or not further substituted in the phenyl
378 ring by one or more other univalent substituents;

379 (ii) By substitution at the 3-position with
380 an alkyl substituent;

381 (iii) By substitution at the nitrogen atom
382 with alkyl or dialkyl groups, or by inclusion of the nitrogen atom
383 in a cyclic structure.



384 (B) The compounds covered in this paragraph (9)
385 include, but are not limited to, any material, compound, mixture
386 or preparation which contains any quantity of a synthetic
387 cathinone found in any of the following compounds, whether or not
388 substituted to any extent, or any of these compounds which contain
389 any synthetic cathinone, or salts, isomers, or salts of isomers,
390 whenever the existence of such salts, isomers or salts of isomers
391 is possible, unless specifically excepted or listed in another
392 schedule:

393 (i) 4-methyl-N-ethylcathinone ("4-MEC");

394 (ii) 4-methyl-alpha-pyrrolidinopropiophenone
395 ("4-MePPP");

396 (iii) Alpha-pyrrolidinopentiophenone
397 (" α -PVP");

398 (iv)
399 1-(1,3-benzodioxol-5-yl)-2-(methylamino)butan-1-one ("butylone");

400 (v) 2-(methylamino)-1-phenylpentan-1-one
401 ("pentedrone");

402 (vi)
403 1-(1,3-benzodioxol-5-yl)-2-(methylamino)pentan-1-one
404 ("pentylone");

405 (vii) 4-fluoro-N-methylcathinone ("4-FMC");

406 (viii) 3-fluoro-N-methylcathinone ("3-FMC");

407 (ix)

408 1-(naphthalen-2-yl)-2-(pyrrolidin-1-yl)pentan-1-one ("naphyrone");



409 (x) Alpha-pyrrolidinobutiophenone ("α-PBP");

410 and

411 (xi)

412 1-(1,3-benzodioxol-5-yl)-2-(ethylamino)-pentan-1-one

413 (N-ethylpentylone, ephylone).

414 (10) (A) Mitragynine; and

415 (B) 7-hydroxymitragynine.

416 **SECTION 2.** This act shall take effect and be in force from

417 and after July 1, 2021.

